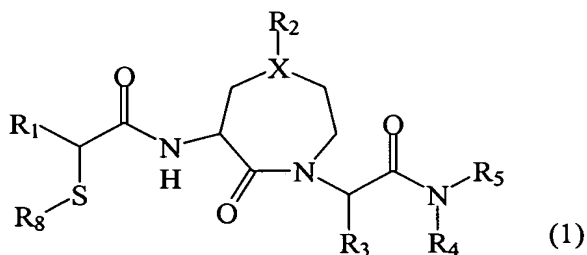


LISTING OF CLAIMS

This listing of claims will replace all prior versions and listings of claims in the application.

1. (Previously presented) A compound of the formula



wherein

R_1 is C_1 - C_6 alkyl, a W -(CH_2) $_m$ - group, or a Q - Z -(CH_2) $_m$ - group wherein W is phthalimido; Z is a bond or is oxy, NR_6 , $C(O)NR_6$, $NR_6C(O)$, $NHC(O)NR_6$, $OC(O)NR_6$, $HNC(O)O$, or SO_2NR_6 ; Q is hydrogen, or a Y -(CH_2) $_n$ - group wherein Y is hydrogen, C_6 - C_{10} aryl, a cyclic or bicyclic, aromatic assemblage of three to nine carbon atoms and from 1 to 3 nitrogen, oxygen, or sulfur atoms, $-C(O)OR_6$, $-N(R_6)_2$, morpholino, piperidino, pyrrolidino, or isoindolyl;

R_2 is C_1 - C_4 alkyl, a $-(CH_2)_p$ -(C_3 - C_9)heteroaryl group, or a $-(CH_2)_p$ - Ar_1 group wherein Ar_1 is phenyl or naphthyl optionally substituted with a substituent selected from the group consisting of halogen, C_1 - C_4 alkyl, $-OR_7$, $-N(R_6)_2$, $SO_2N(R_6)_2$ or $-NO_2$;

R_3 is hydrogen, C_1 - C_6 alkyl, $-CH_2SCH_2NHCOCH_3$, a $-(CH_2)_p$ - A group, a $-(CH_2)_m$ - B group or a $-CH_2$ - D - R_7 group wherein A is C_6 - C_{10} aryl, C_3 - C_9 heteroaryl, or cyclohexyl; B is $-N(R_7)_2$, guanidino, nitroguanidino, $-C(O)OR_6$ or $-C(O)NR_6$; and D is oxy or thio;

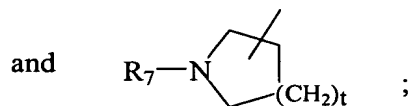
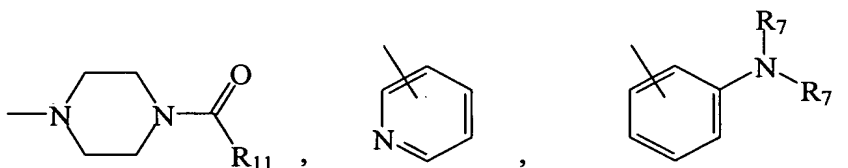
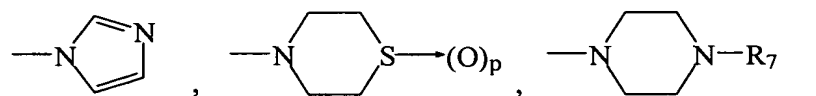
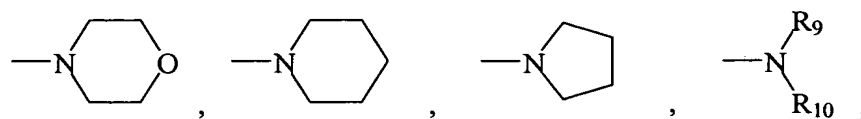
R₄ is hydrogen or a $-(CH_2)_m-S(O)_pX'(R_6)_2$ group;

R₅ is hydrogen, C₁-C₆ alkyl or R₄ and R₅ taken together with the nitrogen atom to which they are attached form piperidino, pyrrolidino, or isoindolyl;

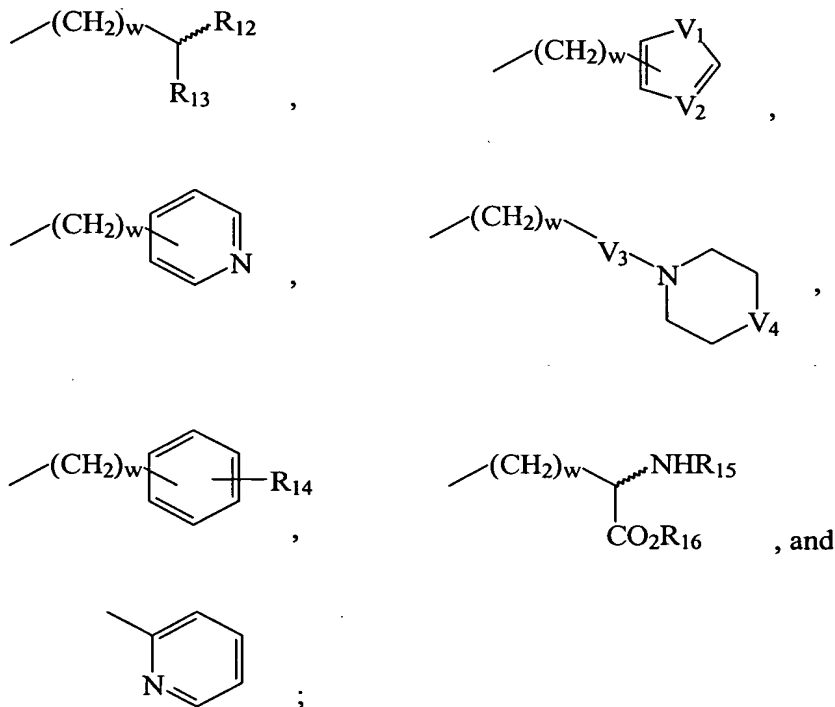
R₆ is hydrogen or C₁-C₆ alkyl;

R₇ is hydrogen, C₁-C₄ alkyl, or a $-(CH_2)_p-Ar_2$ group wherein Ar₂ is phenyl or naphthyl optionally substituted with a substituent selected from the group consisting of halogen, C₁-C₄ alkyl, -OR₇, -N(R₆)₂, SO₂N(R₆)₂ or -NO₂;

R₈ is hydrogen, -C(O)R₇, a -C(O)-(CH₂)_q-K group or a -S-G group, wherein K is selected from the group consisting of



G is selected from the group consisting of



- R₉ and R₁₀ are each independently C₁-C₄ alkyl or a -(CH₂)_p-Ar₂ group;
R₁₁ is -CF₃, C₁-C₁₀ alkyl or a -(CH₂)_p-Ar₂ group;
R₁₂ is hydrogen, C₁-C₆ alkyl, -CH₂CH₂S(O)_pCH₃, or arylalkyl;
R₁₃ is hydrogen, hydroxy, amino, C₁-C₆ alkyl, N-methylamino, N,N-dimethylamino,
-CO₂R₁₇ or -OC(O)R₁₈ wherein R₁₇ is hydrogen, -CH₂O-C(O)C(CH₃)₃,
C₁-C₄ alkyl, a
-(CH₂)_p-Ar₂ group or diphenylmethyl and R₁₈ is hydrogen, C₁-C₆ alkyl
or phenyl;
R₁₄ is 1 or 2 substituents independently chosen from the group consisting of
hydrogen,
C₁-C₄ alkyl, C₁-C₄ alkoxy, or halogen;
R₁₅ is hydrogen, C₁-C₆ alkyl or a -(CH₂)_p-Ar₂ group;

R₁₆ is hydrogen or C₁-C₄ alkyl;
V₁ is O, S, or NH;
V₂ is N or CH;
V₃ is a bond or -C(O)-;
V₄ is -(CH₂)_w-, O, S, NR₇, or NC(O)R₁₁;
X and X' are each independently CH or N;
m is an integer 2-4;
n is zero or an integer 1-4;
p is zero or an integer 1-2;
q is zero or an integer 1-5;
t is an integer 1-2;
w is an integer 1-3; and
w' is zero or an integer 1; or

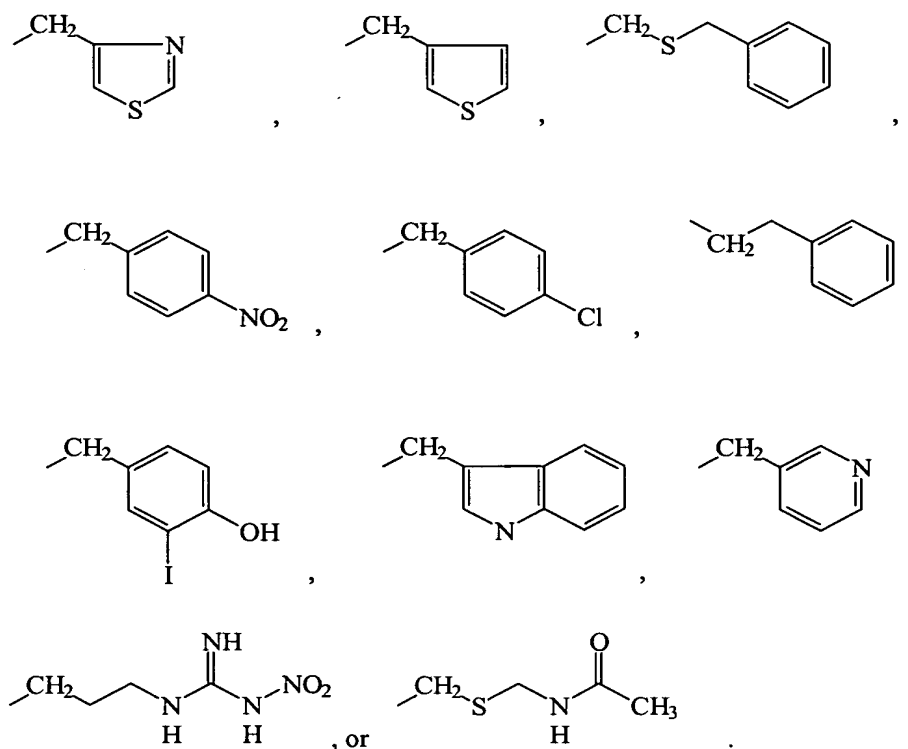
a pharmaceutically acceptable salt, stereoisomer or hydrate thereof.

2. (Original) A compound of claim 1 wherein X is CH.

3. (Original) A compound of claim 2 wherein R₂ is C₁-C₄ alkyl or a -(CH₂)_p-Ar group wherein Ar is phenyl optionally substituted with F, Cl, C₁-C₄ alkyl, -NO₂, -NH₂ or -OR₇; and R₄ is hydrogen.

4. (Previously Presented) A compound of claim 3 wherein R₃ is hydrogen, C₁-C₆ alkyl, phenyl, benzyl, 1-naphthyl, 2-naphthyl, cyclohexylmethyl, 2-hydroxyphenyl, 3-hydroxyphenyl, 4-hydroxyphenyl, 2,3-dihydroxyphenyl, 2,4-dihydroxyphenyl, 3,4-dihydroxyphenyl, 4-methoxyphenyl, 4-ethoxyphenyl, 2-chlorophenyl, 3-chlorophenyl, 4-chlorophenyl, 3,4-dichlorophenyl, 4-bromophenyl, 3,4-dibromophenyl, 4-fluorophenyl, 3,4-difluorophenyl, 3-tolyl, 4-tolyl, 4-ethylphenyl, 4-isopropylphenyl, 3-aminophenyl, 4-aminophenyl, 3,4-diaminophenyl, N-

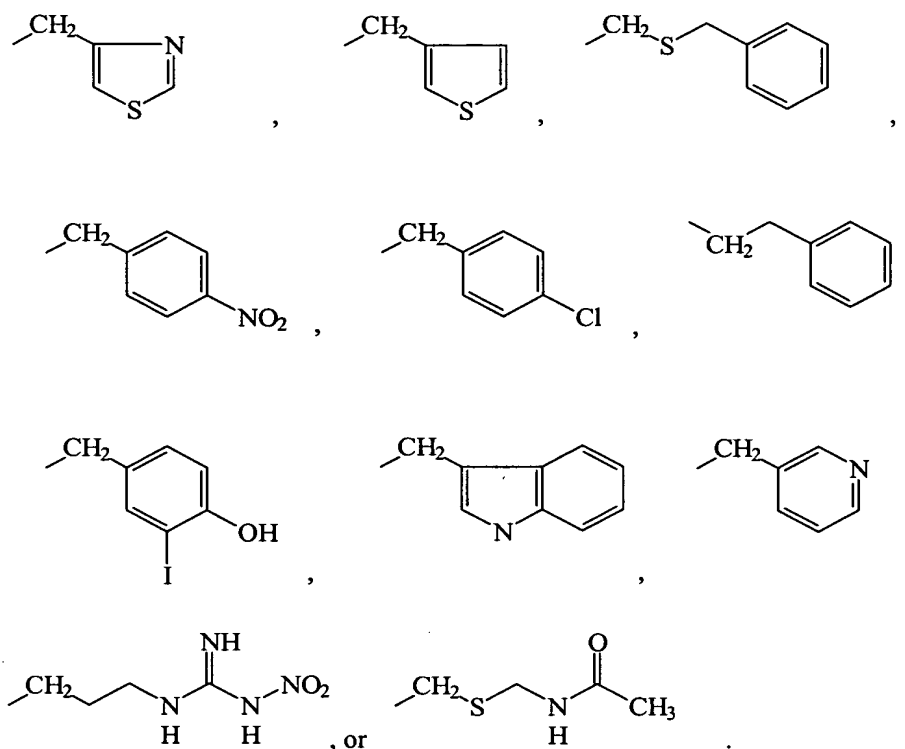
methyl-4-aminophenyl, 2-nitrophenyl, 4-nitrophenyl, 4-aminobenzyl, 4-hydroxybenzyl, 4-methoxybenzyl, 3-chlorobenzyl, 4-fluorobenzyl, 3,4-dichlorobenzyl, 4-bromobenzyl, 4-methylbenzyl, or is a compound of the formula



5. (Original) A compound of claim 4 wherein R_5 is hydrogen, methyl, ethyl, propyl, isopropyl, butyl or isobutyl and R_8 is hydrogen.
6. (Original) A compound of claim 2 wherein R_1 is a $W-(CH_2)_m$ - group.
7. (Original) A compound of claim 3 wherein R_1 is a $W-(CH_2)_m$ - group.
8. (Original) A compound of claim 5 wherein R_1 is a $W-(CH_2)_m$ - group.

9. (Original) A compound of claim 2 wherein R_1 is C_1-C_6 alkyl.
10. (Original) A compound of claim 3 wherein R_1 is C_1-C_6 alkyl.
11. (Original) A compound of claim 5 wherein R_1 is a C_1-C_6 alkyl.
12. (Original) A compound of claim 2 wherein R_1 is a $Q-Z-(CH_2)_m$ -group.
13. (Original) A compound of claim 3 wherein R_1 is a $Q-Z-(CH_2)_m$ -group.
14. (Original) A compound of claim 5 wherein R_1 is a $Q-Z-(CH_2)_m$ -group.
15. (Original) A compound of claim 1 wherein X is N .
16. (Original) A compound of claim 15 wherein R_2 is C_1-C_4 alkyl or a $-(CH_2)_p$ -Ar group wherein Ar is phenyl optionally substituted with F, Cl, C_1-C_4 alkyl, $-NO_2$, $-NH_2$ or $-OR_8$; and R_4 is hydrogen.
17. (Previously Presented) A compound of claim 16 wherein R_3 is hydrogen, C_1-C_6 alkyl, phenyl, benzyl, 1-naphthyl, 2-naphthyl, cyclohexylmethyl, 2-hydroxyphenyl, 3-hydroxyphenyl, 4-hydroxyphenyl, 2,3-dihydroxyphenyl, 2,4-dihydroxyphenyl, 3,4-dihydroxyphenyl, 4-methoxyphenyl, 4-ethoxyphenyl, 2-chlorophenyl, 3-chlorophenyl, 4-chlorophenyl, 3,4-dichlorophenyl, 4-bromophenyl, 3,4-dibromophenyl, 4-fluorophenyl, 3,4-difluorophenyl, 3-tolyl, 4-tolyl, 4-ethylphenyl, 4-isopropylphenyl, 3-aminophenyl, 4-aminophenyl, 3,4-diaminophenyl, N-

methyl-4-aminophenyl, 2-nitrophenyl, 4-nitrophenyl, 4-aminobenzyl, 4-hydroxybenzyl, 4-methoxybenzyl, 3-chlorobenzyl, 4-fluorobenzyl, 3,4-dichlorobenzyl, 4-bromobenzyl, 4-methylbenzyl, or is a compound of the formula



18. (Original) A compound of claim 17 wherein R_5 is hydrogen, methyl, ethyl, propyl, isopropyl, butyl or isobutyl and R_8 is hydrogen.

19. (Original) A compound of claim 16 wherein R_1 is a $W-(CH_2)_m$ -group.

20. (Original) A compound of claim 17 wherein R_1 is a $W-(CH_2)_m$ -group.

21. (Original) A compound of claim 19 wherein R_1 is a $W-(CH_2)_m$ -group.
22. (Original) A compound of claim 16 wherein R_1 is C_1-C_6 alkyl.
23. (Original) A compound of claim 17 wherein R_1 is C_1-C_6 alkyl.
24. (Original) A compound of claim 19 wherein R_1 is a C_1-C_6 alkyl.
25. (Original) A compound of claim 16 wherein R_1 is a $Q-Z-(CH_2)_m$ -group.
26. (Original) A compound of claim 17 wherein R_1 is a $Q-Z-(CH_2)_m$ -group.
27. (Original) A compound of claim 19 wherein R_1 is a $Q-Z-(CH_2)_m$ -group.
28. (Original) A compound of claim 1 wherein X is CH; R_2 is phenyl, methyl or ethyl; R_3 is phenyl, benzyl, cyclohexylmethyl, isopropyl, isobutyl, 3-pyridylmethyl, 4-fluorobenzyl or 4-aminobenzyl; R_4 is hydrogen; R_5 is hydrogen, methyl, ethyl, propyl, isopropyl, butyl or isobutyl and R_8 is hydrogen.
29. (Original) A compound of claim 28 wherein R_1 is a $W-(CH_2)_m$ -group.
30. (Original) A compound of claim 1 wherein X is N; R_2 is phenyl, methyl or ethyl; R_3 is phenyl, benzyl, cyclohexylmethyl, isopropyl, isobutyl, 3-

pyridylmethyl, 4-fluorobenzyl or 4-aminobenzyl; R₄ is hydrogen; R₅ is hydrogen, methyl, ethyl, propyl, isopropyl, butyl or isobutyl and R₈ is hydrogen.

31. (Original) A compound of claim 30 wherein R₁ is a W-(CH₂)_m-group.

32. (Original) A compound of claim 1 wherein said compound is 2*H*-Isoindole-2-hexanamide, *N*-[hexahydro-1-[2-(methylamino)-2-oxo-1-(phenylmethyl)ethyl]-2-oxo-5-phenyl-1*H*-azepin-3-yl]-1,3-dihydro- α -mercapto-1,3-dioxo-, [3*S*-[1(*R**), 3 α , 5 α]]-.

33. (Original) A compound of claim 1 wherein said compound is 2*H*-Isoindole-2-hexanamide, *N*-[hexahydro-1-[2-(methylamino)-2-oxo-1-(phenylmethyl)ethyl]-2-oxo-5-phenyl-1*H*-azepin-3-yl]-1,3-dihydro- α -mercapto-1,3-dioxo-, [3*S*-[1(*R**), 3 α , 5 β]]-.

34. (Original) A compound of claim 1 wherein said compound is 2*H*-Isoindole-2-hexanamide, *N*-[hexahydro-4-[2-(methylamino)-2-oxo-1-(phenylmethyl)ethyl]-5-oxo-1-(phenylmethyl)-1*H*-1,4-diazepin-6-yl]-1,3-dihydro- α -mercapto-1,3-dioxo-, [6*S*-[4(*R**), 6*R**(*R**)]]-.

35. (Original) A pharmaceutical composition comprising a compound of claim 1 and a pharmaceutically acceptable carrier.

36. (Cancelled)

37. (Cancelled)

38. (Original) A method of treating rheumatoid arthritis in a patient in need thereof which comprises administering to the patient an effective matrix metalloproteinase inhibiting amount of a compound of claim 1.

39. (Original) A method of treating osteoarthritis in a patient in need thereof which comprises administering to the patient an effective matrix metalloproteinase inhibiting amount of a compound of claim 1.

40. (Original) A method of treating a chronic inflammatory disorder in a patient in need thereof which comprises administering to the patient an effective matrix metalloproteinase inhibiting amount of a compound of claim 1.

41. (Cancelled)

42. (Cancelled)

43. (Presently amended) A method of ~~claim 42 wherein said cardiovascular disorder is atherosclerosis~~ treating atherosclerosis in a patient in need thereof which comprises administering to the patient an effective matrix metalloproteinase inhibiting amount of a compound of claim 1.

44. (Original) A method of treating corneal ulceration in a patient in need thereof which comprises administering to the patient an effective matrix metalloproteinase inhibiting amount of a compound of claim 1.

45. (Original) A method of treating gingivitis or periodontal disease in a patient in need thereof which comprises administering to the patient an effective matrix metalloproteinase inhibiting amount of a compound of claim 1.

46. (Cancelled)

47. (Original) A method of treating chronic obstructive pulmonary disorder in a patient in need thereof which comprises administering to the patient an effective matrix metalloproteinase inhibiting amount of a compound of claim 1.